

Figure S1. Excitation and emission spectra of compounds **3c**, **3j**, and **3u**.

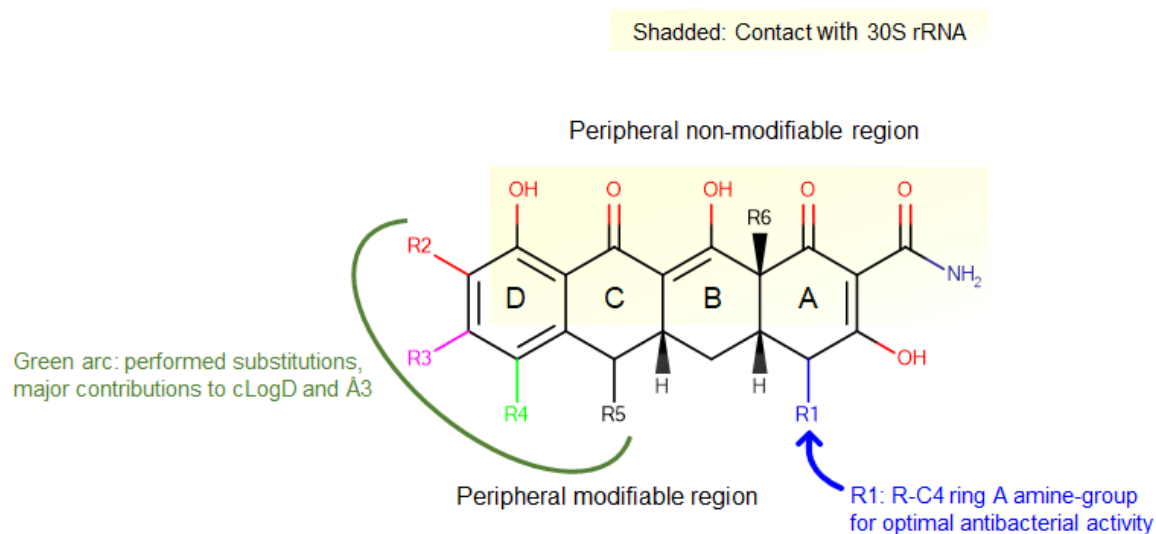


Figure S2. Structure-activity relationship of tetracyclines.

Table S1. List of tetracycline derivatives. Molecules in bold font (**3c**, **3j**, and **3u**) were selected for uptake assays shown in Figure 4 and/or *in vitro* expression assays shown in Figure 5.

Name	n°	MIC (µg/mL)		Fold reduction in MIC	cLogD pH 7.4
		without NV716	with NV716 (10 µM)		
Chlortetracycline	3a	6.25	0.7	9	-4.3
Demeclocycline	3b	6.25	0.4	16	-3.9
Doxycycline	3c	12.5	0.2	63	-2.7
RCT	3d	200	6.25	32	-2.8
RT	3e	400	12.5	32	-2.8
RDOX	3f	400	25	16	-2.7
7-iodo-RDOX	3g	400	6.25	64	-2.5
9-iodo-RDOX	3h	400	12.5	32	-2.8
RDMC	3i	200	6.25	32	-2.4
Col3	3j	400	3.125	128	-1.95
DDOX	3k	400	25	16	-1.8
9-(4-methoxy-phenyl)-RDOX	3l	400	6.25	64	-1.8
9-(3,4-dimethoxyphenyl)-RDOX	3m	400	6.25	64	-1.8
9-(3,4,5-trimethoxyphenyl)- RDOX	3n	400	6.25	64	-1.8
DDMC	3o	250	50	5	-1.9
DCT	3p	400	100	4	-1.6
9-(2-hydroxyphenyl)-RDOX	3q	400	12.5	32	-1.75
9-(3-hydroxyphenyl)-RDOX	3r	400	25	16	-1.75
9-(2-furanyl)-RDOX	3s	400	400	1	-1.4
9-(3,4-methylenedioxyphenyl)- RDOX	3t	400	400	1	-1.2
-	3u	400	400	1	-1.08
(Z)-9-(2-(hydroxymethyl)-hex-1- en-1-yl)-RDOX	3v	400	400	1	-0.9
9-phenyl-RDOX	3w	400	400	1	-0.89
9-tertButyl-Dox-HCl	3x	400	400	1	-0.85
9-(2-thiophenyl)-RDOX	3y	400	400	1	-0.5
9-(2-furanyl)-RDOX	3z	400	400	1	-0.3
9-(2-naphthyl)-RDOX	3aa	400	400	1	0.2
9-(3-benzyloxyphenyl)-RDOX	3ab	400	400	1	0.5